# **TCEQ Interoffice Memorandum**

**To:** Tony Walker

Director, TCEQ Region 4, Dallas/Fort Worth

Alyssa Taylor

Air Section Manager, TCEQ Region 4, Dallas/Fort Worth

From: Manuel Reyna Mark

Toxicology Division, Chief Engineer's Office

**Date:** January 27, 2012

**Subject:** Toxicological Evaluation of Results from an Ambient Air Sample for Volatile

Organic Compounds Collected at Latitude 32.65528, Longitude -97.06654,

Downwind of the Chesapeake Operations, Inc. – West Wing Lease Site in Arlington,

Tarrant County, Texas.

Sample Collected on October 20, 2011, ACL 1111016 (Lab Sample 1111016-001).

## **Key Points**

• Reported concentrations of target volatile organic compounds (VOCs) were either not detected or were detected below levels of short-term health and/or welfare concern.

## **Background**

On October 20, 2011, a Texas Commission on Environmental Quality (TCEQ) Region 4 Air Investigator collected a 30-minute canister sample (Lab Sample 1111016-001) downwind of the Chesapeake Operations, Inc. – West Wing Lease site in Arlington, Tarrant County, Texas (Latitude 32.65528, Longitude -97.06654). The sample was collected in response to a citizen complaint of a burning chemical odor, headache and becoming blind because of a natural gas presence. The investigator did not detect odors or experience health effects during the sampling event. Meteorological conditions measured at the site or nearest stationary ambient air monitoring site indicated that the ambient temperature was 74.5°F with a relative humidity of 29.2%, and winds were from the south at 2.3 miles per hour. The sampling site was greater than 501 feet from the possible emission source (drilling). The nearest location where the public could have access was 101 to 301 feet from the possible emission source. The sample was sent to the TCEQ laboratory in Austin, Texas, and analyzed for a range of VOCs. The list of the target analytes that were evaluated in this review are provided in Attachment A. The VOC concentrations were reported in parts per billion by volume (ppbv) (Attachment B and Table 1). Please note that the available canister technology and analysis method cannot capture and/or analyze for all chemicals.

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#### **Results and Evaluation**

Reported VOC concentrations were compared to TCEQ's short-term health- and/or welfare-based air monitoring comparison values (AMCVs) (Table 1). Short-term AMCVs are guidelines used to evaluate ambient concentrations of a chemical in air and to determine its potential to result in adverse health effects, adverse vegetative effects, or odors. Health AMCVs are set to provide a margin of safety and are set well below levels at which adverse health effects are reported in the scientific literature. If a chemical concentration in ambient air is less than its comparison value, no adverse health effects are expected to occur. If a chemical concentration exceeds its comparison value it does not necessarily mean that adverse effects will occur, but rather that further evaluation is warranted.

All of the 84 VOCs were either not detected or were detected below their respective short-term AMCVs. Exposure to levels of VOCs measured in this sample would not be expected to cause short-term adverse health effects, adverse vegetative effects, or odors.

Please call me at (512) 239-1816 if you have any questions regarding this evaluation.

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#### Attachment A

## **List of Target Analytes for Canister Samples**

ethane ethylene acetylene propane propylene dichlorodifluoromethane methyl chloride isobutane vinyl chloride 1-butene 1.3-butadiene n-butane t-2-butene bromomethane c-2-butene 3-methyl-1-butene

isopentane

trichlorofluoromethane

1-pentene n-pentane isoprene t-2-pentene

1,1-dichloroethylene

c-2-pentene

methylene chloride 2-methyl-2-butene 2,2-dimethylbutane cyclopentene

4-methyl-1-pentene 1.1-dichloroethane cyclopentane 2,3-dimethylbutane 2-methylpentane

3-methylpentane

2-methyl-1-pentene + 1-hexene

n-hexane chloroform t-2-hexene c-2-hexene

1,2-dichloroethane methylcyclopentane 2,4-dimethylpentane 1,1,1-trichloroethane

benzene

carbon tetrachloride

cyclohexane 2-methylhexane 2,3-dimethylpentane 3-methylhexane 1,2-dichloropropane trichloroethylene 2,2,4-trimethylpentane

2-chloropentane

n-heptane

c-1,3-dichloropropylene methylcyclohexane

t-1,3-dichloropropylene 1.1.2-trichloroethane 2,3,4-trimethylpentane

toluene

2-methylheptane 3-methylheptane 1.2-dibromoethane

n-octane

tetrachloroethylene chlorobenzene ethylbenzene m & p-xylene styrene

1,1,2,2-tetrachloroethane

o-xylene n-nonane

isopropylbenzene n-propylbenzene m-ethyltoluene p-ethyltoluene

1,3,5-trimethylbenzene

o-ethyltoluene

1,2,4-trimethylbenzene

n-decane

1,2,3-trimethylbenzene m-diethylbenzene p-diethylbenzene n-undecane

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#### Attachment B

11/16/2011

#### Texas Commission on Environmental Quality

Laboratory and Quality Assurance Section P.O. Box 13087, MC-165 Austin, Texas 78711-3087 (512) 239-1716

#### Laboratory Analysis Results ACL Number: 1111016

ACL Lead: David Manis Region: T04

Project(s): Barnett Shale

Region: T04 Date Received: 11/10/2011

| Facility(ies) Sampled | City      | County  | Facility Type |
|-----------------------|-----------|---------|---------------|
| Chesapeake            | Arlington | Tarrant |               |

#### Laboratory Procedure(s) Performed:

Analysis: AP001VOC

Determination of VOC Canisters by GC/MS Using Modified Method TO-15

Procedure

Prior to analysis, subatmospheric samples are pressurized to twice the collected volume using a sample dilution system. For analysis, a known volume of a sample is directed from the canister into a multitrap cryogenic concentrator. Internal standards are added to the sample stream prior to the trap. The concentrated sample is thermally desorbed and carried onto a GC column for separation. The analytical strategy involves using a GC with dual columns that are coupled to a mass selective detector (MSD) and a flame ionization detector (FID). Mass spectra for individual peaks in the total ion chromatogram are then used for target compound identification and quantitation. The fragmentation pattern is compared with stored spectra taken under similar conditions in order to identify the compound. For any given compound, the intensity of the quantitation fragment is compared with the system response to the fragment for known amounts of the compound. This establishes the compound concentration in the sample. For non-target compound peaks which are at least one-half the height of the internal standard, a library search is performed in an attempt to identify the compound solely upon fracture patterns. These tentatively identified compounds (TIC's) are reported as a sample specific footnote. Accurate quantitation of TICs is not possible. The FID is used for the quantitation of ethane, ethylene, acetylene, propylene and propane and identification is based on matching retention times of standards containing known analytes.

#### Sample(s) Received

Field ID Number: 01202-10202011 Laboratory Sample Number: 1111016-001 Sampled by: Brian Yerkes
Sampling Site: West Wing lease Date & Time Sampled: 10/20/11 16:25:00 Valid Sample: Yes

Comments:

Canister 01202 was used to collect a 30-minute sample using OFC-051.

Please note that this analytical technique is not capable of measuring all compounds which might have adverse health effects. For questions on the analytical procedures please contact the laboratory manager at (512) 239-5853. For an update on the health effects evaluation of these data, please contact the Toxicology Division at (512) 239-1795.

Analyst: Jaydeep Patel

Reviewed By: David Manis (Acting)

Technical Specialist: David Manis

## Laboratory Analysis Results ACL Number: 1111016

Analysis Code: AP001VOC

| Lab ID                                   |      | 1              | 111016-001 |         |               |     |         |
|--|------|----------------|------------|---------|---------------|-----|---------|
| Field ID                                 |      | 01202-10202011 |            |         |               |     |         |
| Canister ID                              |      |                | 01202      |         |               |     |         |
| Analysis Date                            |      |                | 11/15/11   |         |               |     |         |
| Compound                                 | LOD  | Concentration  | SDL        | Flags** | Concentration | SDL | Flags** |
| ethane                                   | 0.50 | 3.6            | 1.0        | T,D1    |               |     | - 1100  |
| ethylene                                 | 0.50 | ND             | 1.0        | T,D1    |               |     |         |
| acetylene                                | 0.50 | ND             | 1.0        | T,D1    |               |     |         |
| propane                                  | 0.50 | 2.0            | 1.0        | L,T,D1  |               |     |         |
| propylene                                | 0.50 | 0.32           | 1.0        | J,T,D1  |               |     |         |
| dichlorodifluoromethane                  | 0.20 | 0.48           | 0.40       | L,D1    |               |     |         |
| methyl chloride                          | 0.20 | 0.56           | 0.40       | L,D1    |               |     |         |
| sobutane                                 | 0.23 | 0.28           | 0.46       | J,D1    |               |     |         |
| vinyl chloride                           | 0.17 | ND             | 0.34       | DI      |               |     |         |
| 1-butene                                 | 0.20 | 0.22           | 0.40       | J,D1    |               |     |         |
| 1,3-butadiene                            | 0.27 | ND             | 0.55       | DI      |               |     |         |
| n-butane                                 | 0.20 | 0.64           | 0.40       | L,D1    |               |     |         |
| 1-2-butene                               | 0.18 | ND             | 0.36       | DI      |               |     |         |
| bromomethane                             | 0.27 | ND             | 0.55       | DI      |               |     |         |
| :-2-butene                               | 0.27 | ND             | 0.55       | DI      |               |     |         |
| 3-methyl-1-butene                        | 0.23 | ND             | 0.46       | DI      |               |     |         |
| sopentane                                | 0.27 | 0.24           | 0.55       | J.D1    | -             |     |         |
| trichlorofluoromethane                   | 0.29 | 0.24           | 0.59       | J,D1    |               |     |         |
| I-pentene                                | 0.27 | ND             | 0.55       | Dl      |               |     |         |
| n-pentane                                | 0.27 | ND             | 0.55       | DI      |               |     |         |
| isoprene                                 | 0.27 | ND             | 0.55       | DI      |               |     |         |
| -2-pentene                               | 0.27 | ND             | 0.55       | DI      | -             |     |         |
| 1,1-dichloroethylene                     | 0.18 | ND             | 0.36       | D1      |               |     |         |
| :-2-pentene                              | 0.25 | ND             | 0.51       | DI      |               |     |         |
| methylene chloride                       | 0.14 | ND             | 0.28       | DI      |               | -   |         |
| 2-methyl-2-butene                        | 0.14 | ND             | 0.46       | DI      |               |     |         |
| 2,2-dimethylbutane                       | 0.23 | ND             | 0.46       | DI      | -             |     |         |
|  |      | ND             |            |         |               |     |         |
| cyclopentene                             | 0.20 |                | 0.40       | D1      |               | -   |         |
| 4-methyl-1-pentene<br>1,1-dichloroethane | 0.22 | ND             | 0.44       | DI      |               |     |         |
|  | 0.19 | ND             | 0.38       | D1      |               |     |         |
| cyclopentane                             | 0.27 | ND             | 0.55       | D1      |               |     |         |
| 2,3-dimethylbutane                       | 0.28 | ND             | 0.57       | DI      |               |     |         |
| 2-methylpentane                          | 0.27 | ND             | 0.55       | D1      |               |     |         |
| 3-methylpentane                          | 0.23 | ND             | 0.46       | Dl      |               |     |         |
| 2-methyl-1-pentene + 1-hexene            | 0.20 | 0.01           | 0.40       | J,D1    |               |     |         |
| n-hexane                                 | 0.20 | ND             | 0.40       | D1      |               |     |         |
| chloroform                               | 0.21 | ND             | 0.42       | D1      |               |     |         |
| -2-hexene                                | 0.27 | ND             | 0.55       | D1      |               |     |         |
| 2-2-hexene                               | 0.27 | ND             | 0.55       | D1      |               |     |         |
| 1,2-dichloroethane                       | 0.27 | ND             | 0.55       | D1      |               |     |         |
| methylcyclopentane                       | 0.27 | ND             | 0.55       | D1      |               |     |         |
| 2,4-dimethylpentane                      | 0.27 | ND             | 0.55       | D1      |               |     |         |
| l,1,1-trichloroethane                    | 0.26 | ND             | 0.53       | D1      |               |     |         |
| benzene                                  | 0.27 | 0.12           | 0.55       | J,D1    |               |     |         |
| arbon tetrachloride                      | 0.27 | 0.10           | 0.55       | J,D1    |               |     |         |
| cyclohexane                              | 0.24 | ND             | 0.48       | D1      |               |     |         |
| 2-methylhexane<br>2,3-dimethylpentane    | 0.27 | ND<br>ND       | 0.55       | D1      |               |     |         |

## **Laboratory Analysis Results**

ACL Number: 1111016 Analysis Code: AP001VOC

| Note: Results are reported in units of pa | erts per billion by vol | ume (ppbv)    |      |         |               |     |         |
|---|-------------------------|---------------|------|---------|---------------|-----|---------|
| ab ID                                     | 1111016-001             |               |      |         |               |     |         |
| Compound                                  | LOD                     | Concentration | SDL  | Flags** | Concentration | SDL | Flags** |
| -methylhexane                             | 0.20                    | ND            | 0.40 | D1      |               |     |         |
| ,2-dichloropropane                        | 0.17                    | ND            | 0.34 | D1      |               |     |         |
| richloroethylene                          | 0.29                    | ND            | 0.59 | D1      |               |     |         |
| ,2,4-trimethylpentane                     | 0.24                    | ND            | 0.48 | D1      |               |     |         |
| -chloropentane                            | 0.27                    | ND            | 0.55 | D1      |               |     |         |
| -heptane                                  | 0.25                    | ND            | 0.51 | D1      |               |     |         |
| -1,3-dichloropropylene                    | 0.20                    | ND            | 0.40 | D1      |               |     |         |
| nethylcyclohexane                         | 0.26                    | ND            | 0.53 | D1      |               |     |         |
| -1,3-dichloropropylene                    | 0.20                    | ND            | 0.40 | D1      |               | -   |         |
| ,1,2-trichloroethane                      | 0.21                    | ND            | 0.42 | D1      |               |     |         |
| ,3,4-trimethylpentane                     | 0.24                    | ND            | 0.48 | D1      |               |     |         |
| oluene                                    | 0.27                    | 0.07          | 0.55 | J,D1    |               |     |         |
| -methylheptane                            | 0.20                    | 0.01          | 0.40 | J,DI    |               |     |         |
| -methylheptane                            | 0.23                    | ND            | 0.46 | D1      |               |     |         |
| ,2-dibromoethane                          | 0.20                    | ND            | 0.40 | D1      |               |     |         |
| -octane                                   | 0.19                    | ND            | 0.38 | D1      |               |     |         |
| etrachloroethylene                        | 0.24                    | ND            | 0.48 | D1      |               |     |         |
| hlorobenzene                              | 0.27                    | ND            | 0.55 | D1      |               |     |         |
| thylbenzene                               | 0.27                    | 0.05          | 0.55 | J,D1    |               |     |         |
| n & p-xylene                              | 0.27                    | 0.03          | 0.55 | J,D1    |               |     |         |
| tyrene                                    | 0.27                    | ND            | 0.55 | D1      |               |     |         |
| ,1,2,2-tetrachloroethane                  | 0.20                    | ND            | 0.40 | D1      |               |     |         |
| -xylene                                   | 0.27                    | ND            | 0.55 | D1      |               |     |         |
| -nonane                                   | 0.22                    | ND            | 0.44 | D1      |               |     |         |
| sopropylbenzene                           | 0.24                    | ND            | 0.48 | D1 -    |               |     |         |
| -propylbenzene                            | 0.27                    | ND            | 0.55 | D1      |               |     |         |
| n-ethyltoluene                            | 0.11                    | ND            | 0.22 | D1      |               |     |         |
| -ethyltoluene                             | 0.16                    | ND            | 0.32 | D1      |               |     |         |
| ,3,5-trimethylbenzene                     | 0.25                    | ND            | 0.51 | D1      |               |     |         |
| -ethyltoluene                             | 0.13                    | ND            | 0.26 | D1      |               |     |         |
| ,2,4-trimethylbenzene                     | 0.27                    | ND            | 0.55 | D1      |               |     |         |
| -decane                                   | 0.27                    | ND            | 0.55 | D1      |               |     |         |
| ,2,3-trimethylbenzene                     | 0.27                    | ND            | 0.55 | D1      |               |     |         |
| n-diethylbenzene                          | 0.27                    | ND            | 0.55 | D1      |               |     |         |
| -diethylbenzene                           | 0.27                    | ND            | 0.55 | D1      |               |     |         |
| -undecane                                 | 0.27                    | ND            | 0.55 | D1      |               |     |         |

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## Laboratory Analysis Results ACL Number: 1111016 Analysis Code: AP001VOC

Note: Results are reported in units of parts per billion by volume ( ppbv)

LOD - Limit of Detection.

ND - not detected

NQ - concentration can not be quantified.

SDL - Sample Detection Limit (LOD adjusted for dilutions).

INV - Invalid.

J - Reported concentration is below SDL.

L - Reported concentration is at or above the SDL and is below the lower limit of quantitation.

E - Reported concentration exceeds the upper limit of instrument calibration.

M - Result modified from previous result.

T- Data was not confirmed by a confirmational analysis. Data is tentatively identified.

\* SDL is equal to LOD

\*\* Quality control flags explanations are listed on the last page of this report.

TCEQ laboratory customer support may be reached at David.Manis@tceq.texas.gov

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### Laboratory Analysis Results ACL Number: 1111016 Analysis Code: AP001VOC

### Quality Control Notes:

D1-sample concentration was calculated using a dilution factor of 4.04

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Table 1. Comparison of Monitored Concentrations in Lab Sample 1111016-001 to TCEQ Short-Term AMCVs

| Lab Sample ID                     | 1111016-001                      |   |                            |                                    |       |                            |
|-----------------------------------|----------------------------------|---|----------------------------|------------------------------------|-------|----------------------------|
| Compound                          | Odor AMCV<br>(ppb <sub>v</sub> ) | Short-Term Health<br>AMCV (ppb <sub>v</sub> ) | LOD<br>(ppb <sub>v</sub> ) | Concentrations (ppb <sub>v</sub> ) | Flags | SDL<br>(ppb <sub>v</sub> ) |
| 1,1,1-Trichloroethane             | 380,000                          | 1,700   | 0.26                       | ND                                 | D1    | 0.53                       |
| 1,1,2,2-Tetrachloroethane         | 7,300                            | 10  | 0.2                        | ND                                 | D1    | 0.4                        |
| 1,1,2-Trichloroethane             | Not Available                    | 100   | 0.21                       | ND                                 | D1    | 0.42                       |
| 1,1-Dichloroethane                | 110,000                          | 1,000   | 0.19                       | ND                                 | D1    | 0.38                       |
| 1,1-Dichloroethylene              | Not Available                    | 180   | 0.18                       | ND                                 | D1    | 0.36                       |
| 1,2,3-Trimethylbenzene            | Not Available                    | 250   | 0.27                       | ND                                 | D1    | 0.55                       |
| 1,2,4-Trimethylbenzene            | Not Available                    | 250   | 0.27                       | ND                                 | D1    | 0.55                       |
| 1,2-Dibromoethane                 | 10,000                           | 0.5   | 0.2                        | ND                                 | D1    | 0.4                        |
| 1,2-Dichloroethane                | 6,000                            | 40  | 0.27                       | ND                                 | D1    | 0.55                       |
| 1,2-Dichloropropane               | 250                              | 100   | 0.17                       | ND                                 | D1    | 0.34                       |
| 1,3,5-Trimethylbenzene            | Not Available                    | 250   | 0.25                       | ND                                 | D1    | 0.51                       |
| 1,3-Butadiene                     | 230                              | 1,700   | 0.27                       | ND                                 | D1    | 0.55                       |
| 1-Butene                          | 360                              | 50,000  | 0.2                        | 0.22                               | J,D1  | 0.4                        |
| 1-Pentene                         | 100                              | 2,600   | 0.27                       | ND                                 | D1    | 0.55                       |
| 2,2,4-Trimethylpentane            | Not Available                    | 750   | 0.24                       | ND                                 | D1    | 0.48                       |
| 2,2-Dimethylbutane (Neohexane)    | Not Available                    | 1,000   | 0.21                       | ND                                 | D1    | 0.42                       |
| 2,3,4-Trimethylpentane            | Not Available                    | 750   | 0.24                       | ND                                 | D1    | 0.48                       |
| 2,3-Dimethylbutane                | Not Available                    | 990   | 0.28                       | ND                                 | D1    | 0.57                       |
| 2,3-Dimethylpentane               | Not Available                    | 850   | 0.26                       | ND                                 | D1    | 0.53                       |
| 2,4-Dimethylpentane               | 290,000                          | 850   | 0.27                       | ND                                 | D1    | 0.55                       |
| 2-Chloropentane (as chloroethane) | Not Available                    | 190   | 0.27                       | ND                                 | D1    | 0.55                       |
| 2-Methyl-1-Pentene +1-Hexene      | 20                               | 500   | 0.2                        | 0.01                               | J,D1  | 0.4                        |
| 2-Methyl-2-Butene                 | 250                              | 500   | 0.23                       | ND                                 | D1    | 0.46                       |
| 2-Methylheptane                   | Not Available                    | 750   | 0.2                        | 0.01                               | J,D1  | 0.4                        |

| Lab Sample ID                   | 1111016-001                      |   |                            |                                    |       |                            |
|---------------------------------|----------------------------------|---|----------------------------|------------------------------------|-------|----------------------------|
| Compound                        | Odor AMCV<br>(ppb <sub>v</sub> ) | Short-Term Health<br>AMCV (ppb <sub>v</sub> ) | LOD<br>(ppb <sub>v</sub> ) | Concentrations (ppb <sub>v</sub> ) | Flags | SDL<br>(ppb <sub>v</sub> ) |
| 2-Methylhexane                  | Not Available                    | 750   | 0.27                       | ND                                 | D1    | 0.55                       |
| 2-Methylpentane (Isohexane)     | 83                               | 1,000   | 0.27                       | ND                                 | D1    | 0.55                       |
| 3-Methyl-1-Butene               | 250                              | 8,000   | 0.23                       | ND                                 | D1    | 0.46                       |
| 3-Methylheptane                 | Not Available                    | 750   | 0.23                       | ND                                 | D1    | 0.46                       |
| 3-Methylhexane                  | Not Available                    | 750   | 0.2                        | ND                                 | D1    | 0.4                        |
| 3-Methylpentane                 | Not Available                    | 1,000   | 0.23                       | ND                                 | D1    | 0.46                       |
| 4-Methyl-1-Pentene (as hexene)  | 20                               | 500   | 0.22                       | ND                                 | D1    | 0.44                       |
| Acetylene                       | 620,000                          | 25,000  | 0.5                        | ND                                 | T,D1  | 1                          |
| Benzene                         | 2,700                            | 180   | 0.27                       | 0.12                               | J,D1  | 0.55                       |
| Bromomethane (methyl bromide)   | 21,000                           | 30  | 0.27                       | ND                                 | D1    | 0.55                       |
| c-1,3-Dichloropropylene         | Not Available                    | 10  | 0.2                        | ND                                 | D1    | 0.4                        |
| c-2-Butene                      | 2,100                            | 15,000  | 0.27                       | ND                                 | D1    | 0.55                       |
| c-2-Hexene                      | Not Available                    | 500   | 0.27                       | ND                                 | D1    | 0.55                       |
| c-2-Pentene                     | Not Available                    | 2,600   | 0.25                       | ND                                 | D1    | 0.51                       |
| Carbon Tetrachloride            | 97,000                           | 20  | 0.27                       | 0.1                                | J,D1  | 0.55                       |
| Chlorobenzene (phenyl chloride) | 210                              | 100   | 0.27                       | ND                                 | D1    | 0.55                       |
| Chloroform (trichloromethane)   | 85,000                           | 20  | 0.21                       | ND                                 | D1    | 0.42                       |
| Cyclohexane                     | 420                              | 1,000   | 0.24                       | ND                                 | D1    | 0.48                       |
| Cyclopentane                    | Not Available                    | 1,200   | 0.27                       | ND                                 | D1    | 0.55                       |
| Cyclopentene                    | Not Available                    | 2,900   | 0.2                        | ND                                 | D1    | 0.4                        |
| Dichlorodifluoromethane         | Not Available                    | 10,000  | 0.2                        | 0.48                               | L,D1  | 0.4                        |
| Ethane                          | 180,000                          | Simple Asphyxiant*                            | 0.5                        | 3.6                                | T,D1  | 1                          |
| Ethylbenzene                    | 170                              | 20,000  | 0.27                       | 0.05                               | J,D1  | 0.55                       |
| Ethylene                        | 270,000                          | 500,000                                       | 0.5                        | ND                                 | T,D1  | 1                          |
| Isobutane                       | 2,040                            | 8,000   | 0.23                       | 0.28                               | J,D1  | 0.46                       |
| Isopentane (2-methylbutane)     | 1,300                            | 68,000  | 0.27                       | 0.24                               | J,D1  | 0.55                       |

| Lab Sample ID                        | 1111016-001                      |   |                            |                                       |        |                            |
|--------------------------------------|----------------------------------|---|----------------------------|---------------------------------------|--------|----------------------------|
| Compound                             | Odor AMCV<br>(ppb <sub>v</sub> ) | Short-Term Health<br>AMCV (ppb <sub>v</sub> ) | LOD<br>(ppb <sub>v</sub> ) | Concentrations<br>(ppb <sub>v</sub> ) | Flags  | SDL<br>(ppb <sub>v</sub> ) |
| Isoprene                             | 5                                | 20  | 0.27                       | ND                                    | D1     | 0.55                       |
| Isopropylbenzene (cumene)            | 100                              | 500   | 0.24                       | ND                                    | D1     | 0.48                       |
| m & p-Xylene (as mixed isomers)      | 80                               | 1,700   | 0.27                       | 0.03                                  | J,D1   | 0.55                       |
| m-Diethylbenzene                     | 70                               | 460   | 0.27                       | ND                                    | D1     | 0.55                       |
| Methyl Chloride (chloromethane)      | Not Available                    | 500   | 0.2                        | 0.56                                  | L,D1   | 0.4                        |
| Methylcyclohexane                    | 150                              | 4,000   | 0.26                       | ND                                    | D1     | 0.53                       |
| Methylcyclopentane                   | 1,700                            | 750   | 0.27                       | ND                                    | D1     | 0.55                       |
| Methylene Chloride (dichloromethane) | 160,000                          | 3,500   | 0.14                       | ND                                    | D1     | 0.28                       |
| m-Ethyltoluene                       | 18                               | 250   | 0.11                       | ND                                    | D1     | 0.22                       |
| n-Butane                             | 1,200,000                        | 8,000   | 0.2                        | 0.64                                  | L,D1   | 0.4                        |
| n-Decane                             | 620                              | 1,750   | 0.27                       | ND                                    | D1     | 0.55                       |
| n-Heptane                            | 670                              | 850   | 0.25                       | ND                                    | D1     | 0.51                       |
| n-Hexane                             | 1,500                            | 1,800   | 0.2                        | ND                                    | D1     | 0.4                        |
| n-Nonane                             | 2,200                            | 2,000   | 0.22                       | ND                                    | D1     | 0.44                       |
| n-Octane                             | 1,700                            | 750   | 0.19                       | ND                                    | D1     | 0.38                       |
| n-Pentane                            | 1,400                            | 68,000  | 0.27                       | ND                                    | D1     | 0.55                       |
| n-Propylbenzene                      | 3.8                              | 250   | 0.27                       | ND                                    | D1     | 0.55                       |
| n-Undecane                           | Not Available                    | 550   | 0.27                       | ND                                    | D1     | 0.55                       |
| o-Ethyltoluene                       | Not Available                    | 250   | 0.13                       | ND                                    | D1     | 0.26                       |
| o-Xylene                             | 380                              | 1,700   | 0.27                       | ND                                    | D1     | 0.55                       |
| p-Diethylbenzene                     | 0.39                             | 460   | 0.27                       | ND                                    | D1     | 0.55                       |
| p-Ethyltoluene                       | 8.3                              | 250   | 0.16                       | ND                                    | D1     | 0.32                       |
| Propane                              | 1,500,000                        | Simple Asphyxiant*                            | 0.5                        | 2                                     | L,T,D1 | 1                          |
| Propylene                            | 13,000                           | Simple Asphyxiant*                            | 0.5                        | 0.32                                  | J,T,D1 | 1                          |
| Styrene                              | 25                               | 5,100   | 0.27                       | ND                                    | D1     | 0.55                       |
| t-1,3-Dichloropropylene              | Not Available                    | 10  | 0.2                        | ND                                    | D1     | 0.4                        |

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| Lab Sample ID          | 1111016-001                      | 1111016-001                                   |                            |                                    |       |                            |  |  |
|------------------------|----------------------------------|---|----------------------------|------------------------------------|-------|----------------------------|--|--|
| Compound               | Odor AMCV<br>(ppb <sub>v</sub> ) | Short-Term Health<br>AMCV (ppb <sub>v</sub> ) | LOD<br>(ppb <sub>v</sub> ) | Concentrations (ppb <sub>v</sub> ) | Flags | SDL<br>(ppb <sub>v</sub> ) |  |  |
| t-2-Butene             | 2,100                            | 15,000  | 0.18                       | ND                                 | D1    | 0.36                       |  |  |
| t-2-Hexene             | Not Available                    | 500   | 0.27                       | ND                                 | D1    | 0.55                       |  |  |
| t-2-Pentene            | Not Available                    | 2,600   | 0.27                       | ND                                 | D1    | 0.55                       |  |  |
| Tetrachloroethylene    | 770                              | 1,000   | 0.24                       | ND                                 | D1    | 0.48                       |  |  |
| Toluene                | 170                              | 4,000   | 0.27                       | 0.07                               | J,D1  | 0.55                       |  |  |
| Trichloroethylene      | 3,900                            | 100   | 0.29                       | ND                                 | D1    | 0.59                       |  |  |
| Trichlorofluoromethane | 5,000                            | 10,000  | 0.29                       | 0.24                               | J,D1  | 0.59                       |  |  |
| Vinyl Chloride         | Not Available                    | 26,000  | 0.17                       | ND                                 | D1    | 0.34                       |  |  |

<sup>\*</sup>A simple asphyxiant displaces air, lowering the partial pressure of oxygen and causing hypoxia at sufficiently high concentrations. ppbv - Parts per billion by volume.

ND - Not detected.

NQ - Concentration cannot be quantified.

LOD - Limit of detection.

SDL - Sample Detection Limit (LOD adjusted for dilutions).

INV - Invalid.

- J Reported concentration is below SDL.
- L Reported concentration is at or above the SDL and is below the lower limit of quantitation.
- E Reported concentration exceeds the upper limit of instrument calibration.
- M Result modified from previous result.
- T Data was not confirmed by a confirmational analysis. Data is tentatively identified.
- C Sample received with missing or broken custody seal.
- D1 Sample concentration was calculated using a dilution factor of 4.04.

**Table 2. TCEQ Long-Term Air Monitoring Comparison Values (AMCVs)** 

Please Note: The long-term AMCVs are provided for informational purposes only because it is scientifically inappropriate to compare short-term monitored values to the long-term AMCV.

| Compound                          | Long-Term Health AMCV (ppb <sub>v</sub> ) | Compound                             | Long-Term Health<br>AMCV (ppb <sub>v</sub> ) |
|-----------------------------------|---|--------------------------------------|--|
| 1,1,1-Trichloroethane             | 940                                       | Cyclopentane                         | 120  |
| 1,1,2,2-Tetrachloroethane         | 1   | Cyclopentene                         | 290  |
| 1,1,2-Trichloroethane             | 10  | Dichlorodifluoromethane              | 1,000  |
| 1,1-Dichloroethane                | 100                                       | Ethane                               | Simple Asphyxiant*                           |
| 1,1-Dichloroethylene              | 86  | Ethylbenzene                         | 450  |
| 1,2,3-Trimethylbenzene            | 25  | Ethylene**                           | 5,300  |
| 1,2,4-Trimethylbenzene            | 25  | Isobutane                            | 800  |
| 1,2-Dibromoethane                 | 0.05                                      | Isopentane (2-methylbutane)          | 8,000  |
| 1,2-Dichloroethane                | 1   | Isoprene                             | 2  |
| 1,2-Dichloropropane               | 10  | Isopropylbenzene (cumene)            | 50   |
| 1,3,5-Trimethylbenzene            | 25  | m & p-Xylene (as mixed isomers)      | 140  |
| 1,3-Butadiene                     | 9.1                                       | m-Diethylbenzene                     | 46   |
| 1-Butene                          | Not Available                             | Methyl Chloride (chloromethane)      | 50   |
| 1-Pentene                         | Not Available                             | Methylcyclohexane                    | 400  |
| 2,2,4-Trimethylpentane            | 75  | Methylcyclopentane                   | 75   |
| 2,2-Dimethylbutane (Neohexane)    | 100                                       | Methylene Chloride (dichloromethane) | 100  |
| 2,3,4-Trimethylpentane            | 75  | m-Ethyltoluene                       | 25   |
| 2,3-Dimethylbutane                | 99  | n-Butane                             | 800  |
| 2,3-Dimethylpentane               | 85  | n-Decane                             | 175  |
| 2,4-Dimethylpentane               | 85  | n-Heptane                            | 85   |
| 2-Chloropentane (as chloroethane) | 19  | n-Hexane                             | 190  |
| 2-Methyl-1-Pentene +1-Hexene      | 50  | n-Nonane                             | 200  |

| Compound                        | Long-Term Health AMCV (ppb <sub>v</sub> ) | Compound                | Long-Term Health<br>AMCV (ppb <sub>v</sub> ) |
|---------------------------------|---|-------------------------|--|
| 2-Methyl-2-Butene               | 50  | n-Octane                | 75   |
| 2-Methylheptane                 | 75  | n-Pentane               | 8,000  |
| 2-Methylhexane                  | 75  | n-Propylbenzene         | 25   |
| 2-Methylpentane (Isohexane)     | 100                                       | n-Undecane              | 55   |
| 3-Methyl-1-Butene               | 800                                       | o-Ethyltoluene          | 25   |
| 3-Methylheptane                 | 75  | o-Xylene                | 140  |
| 3-Methylhexane                  | 75  | p-Diethylbenzene        | 46   |
| 3-Methylpentane                 | 100                                       | p-Ethyltoluene          | 25   |
| 4-Methyl-1-Pentene (as hexene)  | 50  | Propane                 | Simple Asphyxiant*                           |
| Acetylene                       | 2,500                                     | Propylene               | Simple Asphyxiant*                           |
| Benzene                         | 1.4                                       | Styrene                 | 110  |
| Bromomethane (methyl bromide)   | 3   | t-1,3-Dichloropropylene | 1  |
| c-1,3-Dichloropropylene         | 1   | t-2-Butene              | Not Available                                |
| c-2-Butene                      | Not Available                             | t-2-Hexene              | 50   |
| c-2-Hexene                      | 50  | t-2-Pentene             | Not Available                                |
| c-2-Pentene                     | Not Available                             | Tetrachloroethylene***  | 3.8  |
| Carbon Tetrachloride            | 2   | Toluene                 | 1,100  |
| Chlorobenzene (phenyl chloride) | 10  | Trichloroethylene       | 10   |
| Chloroform (trichloromethane)   | 2   | Trichlorofluoromethane  | 1,000  |
| Cyclohexane                     | 100                                       | Vinyl Chloride          | 0.45   |

<sup>\*</sup>A simple asphyxiant displaces air, lowering the partial pressure of oxygen and causing hypoxia at sufficiently high concentrations.

<sup>\*\*</sup>Long-term vegetation AMCV for Ethylene is 30 ppb.

<sup>\*\*\*</sup>Long-term vegetation AMCV for Tetrachloroethylene is 12 ppb.